

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092518\
 Data File : BF109288.D
 Acq On : 25 Sep 2018 16:25
 Operator : JU/SJ
 Sample : J4774-05 2X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-03-092418

Manual Integrations
 APPROVED

Sohil
 9/26/2018 3:41:50 PM

Quant Time: Sep 26 02:30:10 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 22 13:32:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.71	152	121802	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	474854	20.00	ng	0.00
38) Acenaphthene-d10	9.74	164	222352	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	361736	20.00	ng	0.00
75) Chrysene-d12	13.84	240	307230	20.00	ng	0.00
86) Perylene-d12	15.25	264	260162	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.31	112	63340	8.57	ng	0.00
7) Phenol-d6	6.34	99	82312	8.72	ng	-0.02
23) Nitrobenzene-d5	7.26	82	48195	5.99	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	15730	7.18	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	102182	6.93	ng	-0.01
78) Terphenyl-d14	12.80	244	80043	6.39	ng	0.00
Target Compounds						
70) Phenanthrene	11.24	178	63324	3.506	ng	Qvalue 96
74) Fluoranthene	12.42	202	81011	4.197	ng	97
77) Pyrene	12.65	202	74379	3.378	ng	97
87) Benzo(b)fluoranthene	14.85	252	36768m	2.572	ng	
89) Benzo(a)pyrene	15.19	252	26027	2.020	ng	# 91

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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